40

Claims

1. A compound corresponding to formula (I)

$$R^{1} - X - C - R^{3}$$
 N
 R^{4}
 R^{5}

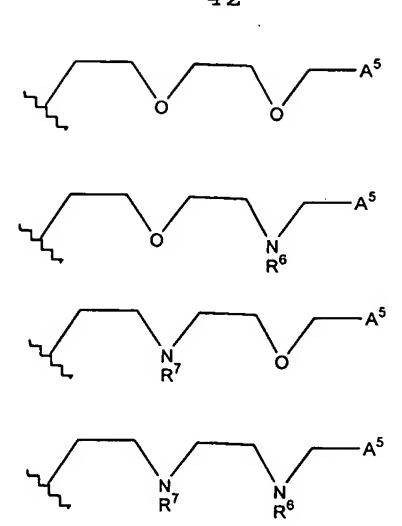
10 in which

5

- R¹ is a functional group capable of reacting with the functions present on proteins, antibodies or on mineral or organic materials;
- X represents a single bond or a hydrocarbon-based chain consisting of at least one group chosen from alkylene groups and alkenylene groups optionally comprising at least one hetero atom, and from arylene groups;
- R^2 is a group A^2 that is anionic at neutral pH or an alkylene or alkenylene group containing from 1 to 4 carbon atoms and bearing at least one such group A^2 , said alkylene or alkenylene group optionally comprising at least one hetero atom in the chain;
- R³ represents H or an alkylene or alkenylene group containing from 1 to 5 carbon atoms and optionally containing at least one hetero atom in the chain, said group optionally bearing at least one group A³ that is anionic at neutral pH;
- R⁴ is chosen from the groups corresponding to the formula $-(C)_n-C-Z^1-C-C-Z^2-C-A^4$ in which n is equal to 1 or 2, 30 Z¹ and Z² represent, independently of each other, a hetero atom chosen from O and N, at least one being a nitrogen atom forming part of an aromatic heterocycle with the two carbon atoms surrounding it, and A⁴ is a group that is anionic at neutral pH, in which the atom bearing the anionic charge is in the γ position relative to Z²;
 - $^ R^5$ is chosen from the groups defined for R^4 or from groups corresponding to the formula $-C-C-E^1-C-C-E^2-C-A^5$ in which E^1 and E^2 represent, independently of each other, a

hetero atom chosen from O and N, and A^5 is a group that is anionic at neutral pH, in which the atom bearing the anionic charge is in the γ position relative to E^2 .

- 2. The compound as claimed in claim 1, characterized in that the substituent R¹ is chosen from amino, thio, cyano, isocyano, acridinyl, hydrazino, haloacetate, anhydride, triazo, carbonyl, nitrobenzoyl, sulfonyl, thionyl, halide, epoxide, aldehyde, imidazole, hydroxyphenyl, mercapto, N-succinimidyl ester, N-sulfosuccinimidyl ester, maleimido, hydroxyl, carboxyl, thiocyano, and isothiocyano groups.
 - 3. The compound as claimed in claim 1, characterized in that the substituent R^2 is a group A^2 that is anionic at neutral pH.
- 4. The compound as claimed in claim 1, characterized in that the substituent R^3 is H or a C_1 to C_3 alkyl.
 - 5. The compound as claimed in claim 1, characterized in that the groups Z^1 and Z^2 of R^4 form part of an aromatic heterocyclic group.
- 20 6. The compound as claimed in claim 1, characterized in that n is equal to 1.
- 7. The compound as claimed in claim 1, characterized in that one of the segments $-C-Z^1-C-$ or $-C-Z^2-C-$ forms part of a heterocyclic group chosen from pyridyl, pyrimidinyl, quinolyl and isoquinolyl groups.
- 8. The compound as claimed in claim 1, characterized in that the segment $-C-Z^1-C-C-Z^2-C-$ is chosen from 2,2'-bipyridinyl, 1,10-phenanthrolinyl, 2,2'-bisquinolyl, 2,2'-bisisoquinolyl and 2,2'-bipyrimidinyl groups, said groups possibly bearing alkyl or alkoxy substituents on at least one carbon atom of a heterocycle.
 - 9. The compound as claimed in claim 1, characterized in that ${\bf R}^5$ is chosen from the following groups:



in which R^6 and R^7 represent alkyl chains containing from 1 to 5 carbon atoms and optionally containing one or more hetero atoms.

- 10. The compound as claimed in claim 1, characterized in that R^4 and R^5 are identical.
- 11. The compound as claimed in claim 1, characterized in that the groups A^2 , A^3 , A^4 or A^5 that are anionic at neutral pH are chosen, independently of each other, from $-CO_2H$, $-SO_3H$, -P(O)(OR)OH, -P(O)R(OH) and $-P(O)(OH)_2$ groups in which R is an alkyl group or an aryl group.
- 12. The compound as claimed in claim 1, characterized in that it is in cationic form, the nitrogen bearing the substituents R^4 and R^5 , and also possibly the hetero atoms Z^1 , Z^2 , E^1 and E^2 , being in protonated form.
 - 13. The compound as claimed in claim 1, characterized in that it is in anionic form, the various groups A^i being in the form of salts.
- 14. The compound as claimed in claim 1, characterized in that it is in zwitterionic form, the nitrogen bearing the substituents R^4 and R^5 , and also possibly the hetero atoms Z^1 , Z^2 , E^1 and E^2 , being in protonated form, and the various groups A^i being in the form of salts.
- 15. The compound as claimed in claim 1, characterized in that X is an arylene group comprising one or more fused or unfused aromatic nuclei, said nucleus (nuclei) optionally bearing one or more aliphatic hydrocarbon-based groups.

- 16. The compound as claimed in claim 1, characterized in that the group X is an alkylene or alkenylene group containing from 1 to 10 carbon atoms.
- 17. The compound as claimed in claim 1, characterized in that the group X is an arylene group containing from 5 to 10 carbon atoms.
- 18. A process for preparing a lanthanide complex, characterized in that it consists in reacting a compound (I) as claimed in any one of claims 1 to 17 with a compound giving a lanthanide cation.
 - 19. The process as claimed in claim 18, characterized in that the compound giving a lanthanide cation is chosen from lanthanide halide hydrates, lanthanide nitrate hydrates, lanthanide carbonates and lanthanide triflates.
- 20. The process as claimed in claim 18, characterized in that the reaction is performed in solution in a solvent chosen from water, methanol, ethanol and acetonitrile.
- 21. The process as claimed in claim 18, characterized in that compound (I) is reacted with the lanthanide ion precursor in a mixture of methanol and water at a pH ranging from 3 to 5, for a time of between 10 minutes and 24 hours, at a temperature of between 25°C and 80°C, and the pH of the solution is then adjusted to 7.0 and the methanol is evaporated off.
- 22. A complex obtained via a process as claimed in claim 18, consisting of a lanthanide ion Ln complexed with a ligand corresponding to formula (I).
 - 23. The complex as claimed in claim 22, characterized in that the lanthanide ion is chosen from europium, terbium, samarium, dysprosium, erbium, ytterbium, neodymium and gadolinium ions.
- 24. The complex as claimed in claim 22, characterized in that the substituent R^4 of the compound (I) is $-C-C-Z^1-C-C-Z^2-C-A^4$, the 3 chelate rings being formed between the lanthanide cation and, respectively:
 - the N atom bearing R^4 and R^5 , Z^1 and the carbon atoms that separate them;

- Z^1 , Z^2 and the two carbon atoms that separate them;
- the end segment Z^2-C-A^4 .
- 25. The complex as claimed in claim 24, characterized in that the substituent R^5 is of the same type as the substituent R^4 .
 - 26. The complex as claimed in claim 24, characterized in that the substituent R^5 is of the type $-C-C-E^1-C-C-E^2-C-A^5$, three 5-membered chelate rings being formed between the lanthanide cation and, respectively:
- 10 the N atom bearing R⁴ and R⁵, E¹ and the two carbon atoms that separate them;
 - E^1 , E^2 and the two carbon atoms that separate them;
 - the end segment E^2-C-A^5 .
- 27. A process for the quantitative or qualitative analysis of a compound, characterized in that it consists in covalently bonding to said compound a marker consisting of a complex as claimed in one of claims 25 to 29, and in detecting or quantifying the presence of the marked compound by means of the luminescence properties of the marker.
- 28. The process as claimed in claim 27, characterized in that the complex is a europium, terbium, samarium or dysprosium complex.
- 29. The process as claimed in claim 27, characterized in that the substituent R¹ of the complex is chosen from 25 amino, thio and carboxyl groups or from maleimido, N-succinimidyl ester and isothiocyano groups.
 - 30. A relaxation agent for nuclear magnetic resonance, consisting of a complex as claimed in one of claims 22 to 26.
- 31. The relaxation agent as claimed in claim 30, characterized in that it consists of a gadolinium, europium or dysprosium complex.
- 32. The relaxation agent as claimed in claim 30, characterized in that it consists of a complex in which the substituent R^1 is chosen from amino, thio and carboxyl groups or from maleimido, N-succinimidyl ester and isothiocyano groups.